

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN050719\
 Data File : BN005511.D
 Acq On : 07 May 2019 06:14
 Operator : JU/SJ
 Sample : K2403-21
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A5134

Manual Integrations
 APPROVED

mohammad
 5/8/2019 12:01:26 AM

Quant Time: May 07 08:07:52 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN050719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 07 07:04:57 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	2097	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.35	136	8508	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.23	164	5550	0.40	ng/ul	-0.02
10) Phenanthrene-d10	16.98	188	13423m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.22	240	13635m	0.40	ng/ul	0.00
20) Perylene-d12	23.43	264	12119m	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.97	152	3428	0.26	ng/ul	-0.02
14) Fluoranthene-d10	19.03	212	10200m	0.24	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.40	128	1722	0.076	ng/ul#	92
5) 2-Methylnaphthalene	12.05	142	1318	0.086	ng/ul	94
7) Acenaphthylene	13.94	152	2350	0.084	ng/ul	95
9) Fluorene	15.28	166	3262	0.141	ng/ul	98
12) Phenanthrene	17.02	178	30014	0.821	ng/ul	99
13) Anthracene	17.11	178	2592m	0.071	ng/ul	
15) Fluoranthene	19.06	202	30652m	0.606	ng/ul	
17) Pyrene	19.42	202	28756m	0.571	ng/ul	
18) Benzo(a)anthracene	21.20	228	11393	0.242	ng/ul	96
19) Chrysene	21.26	228	18427	0.319	ng/ul	97
21) Benzo(b)fluoranthene	22.77	252	25151m	0.690	ng/ul	
22) Benzo(k)fluoranthene	22.81	252	8769m	0.192	ng/ul	
23) Benzo(a)pyrene	23.33	252	14498m	0.356	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.63	276	10841m	0.238	ng/ul	
25) Dibenzo(a,h)anthracene	25.63	278	2296m	0.061	ng/ul	
26) Benzo(a,h,i)perylene	26.30	276	8282m	0.213	ng/ul	

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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